IN THE CLAIMS

The following is an updated listing of the claims in the application with claims 1, 3, 4, 6-11 and 12 shown as currently amended.

LISTING OF CLAIMS

(Currently amended) An indene derivative A compound of formula
 or a pharmaceutically acceptable salt thereof:

wherein.

 R_1 is C_{1-6} alkyl, C_{1-6} alkenyl, or C_{3-6} cycloalkyl, each of which is unsubstituted or substituted substituted with one or more phenyl groups;

being each unsubstituted or substituted substituted with one or more substituents selected from the group

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consisting of halogen, CN, NH_2 , NO_2 , OR^a , phenyloxy, $C_{1.6}$ alkyl, and $C_{3.6}$ cycloalkyl; and

 R_4 , R_5 , R_6 , and R_7 are each independently H, OH, OSO₂CH₃, O(CH₂)_mR^e, CH₂R^f, OCOCH₂OR^g, OCH₂CH₂OR^g, OCH₂CH=CHR^g, eppyridine 2 yloxy₁ or R_5 and R_6 together form OCH₂O;

in which R^a is H, e_F C_{1-6} alkyl, or C_{3-6} cycloalkyl, C_{1-6} alkyl and C_{3-6} cycloalkyl being each unsubstituted or substituted substituted with one or more halogens;

 R^{b} and R^{c} are each independently H, $C_{\text{1-6}}$ alkyl, or $C_{\text{3-6}}$ cycloalkyl;

Rd is O. S. or NRa;

Re is H, halogen, C₃₋₆ cycloalkyl, naphthyl,

phenyl being unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, CF₃, and COOR^a:

$$R^f$$
 is $OCH_2CH_2R^g$ or $-\frac{1}{2}N$ R^d ;

R^g is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, and OR^a: and

m is an integer in the range of 1 to 5.

(currently amended) The compound of claim 1, wherein R₁ is C_{1.6} alkyl, which is unsubstituted or substituted substituted with a phenyl group; R₂ is H, CN, CO₂R^a, CH₂CO₂R^a, CONR^bR^c, or phenyl; R₃ is C_{1.6} alkyl, C_{3.6}

3. (currently amended) The compound of claim 2, wherein R_1 is CH_3 ; R_2 is H, CN, CO_2R^a , or $CONR^bR^c$; R_3 is C_{1-6} alkyl, er phenyl, or R_1 , phenyl being unsubstituted or

substitutied substituted with one or more halogens or C_{1-6} alkyl groups; and R_5 and R_6 are each independently $O(CH_2)_m R^e$ or $CH_2 R^f$, or together form $OCH_2 O$.

- 4. (currently amended) The \underline{A} compound of claim 1, which is selected from the group consisting of:
- 1) 6-methoxy-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 2) 1-(*trans*-isopropylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 3) 1-(*trans*-benzylimino-*N*-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 4) 1-(trans-ethylimino-N-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 5) 6-methoxy-1-(*trans*-phenylpropylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 6) 6-methoxy-1-(*trans*-(2-methylbutenylimino)-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 7) 1-(trans-isobutylimino-N-oxy)-6-methoxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 8) 1-(trans-methylimino-N-oxy)-6-(2-morphorline-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 9) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 10) 1-(*trans*-methylimino-*N*-oxy)-6-phenetyloxy-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 11) 3-furan-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 12) 6-hydroxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 13) 1-(cis-methylimino-N-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 14) 3-(trans-methylimino-N-oxy)-1-phenyl-3H-indene-5-ol
- 15) 1-(trans-methylimino-N-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- 16) 1-(cis-methylimino-N-oxy)-3-phenyl-6-(5-phenylpentyloxy)-1H-indene-2-carboxylate ethyl ester
- $17) \ 6-[2-(4-chlorophenoxy)acetoxy]-1-({\it trans}-methylimino-N-oxy})-3-phenyl-1 H-indene-2-carboxylate ethyl ester$

18) 6-[2-(4-chlorophenoxy)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester

- 19) 1-(*trans*-methylimino-*N*-oxy)-6-(naphthalene-2-ylmethoxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 20) methyl-[3-phenyl-6-(3-phenylpropoxy)indene-1-yllidene]amine-N-oxide
- 21) 1-(*trans*-methylimino-*N*-oxy)-6-[2-(5-methyl-2-phenylthiazol-4-yl)ethoxy]-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 22) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 23) 6-[2-(4-hydroxyphenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 24) 6-(2-adaman-1-ylethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 25) 6-(2-cyclohexylethoxy)-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-
- 2-carboxylate ethyl ester
 26) 1-(trans-methylimino-N-oxy)-3-phenyl-6-(3-phenylprophenoxy)-1H-indene2-carboxylate ethyl ester
- 27) 6-[2-(2-fluorophenyl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 28) 6-[2-(3-fluorophenyl)ethoxy]-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- Indene-2-carboxylate etnyl ester 29) 6-[2-(4-fluorophenyl)ethoxy]-1-(trans-methylimino-N-oxy)-3-phenyl-1Hindene-2-carboxylate etnyl ester
- 30) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-[2-(3-trifluoromethylphenyl)ethoxyl-1H-indene-2-carboxylate ethyl ester
- 31) 6-(4-methoxycarbonylbenzyloxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 32) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-
- carboxylate ethyl amide
 33) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1Hindene-2-carboxylate ethyl ester
- 34) 6-[2-(cyclohexylmethylamino)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
 - 35) 3-(2-fluorophenyl)-6-methoxy-1-(*trans*-methylimino-*N*-oxy)-1H-indene-2-carboxylate ethyl ester
 - 36) 1-(trans-methylimino-N-oxy)-6-[2-(4-methylpiperazine-1-yl)ethoxy]-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 37) (2,3-diphenyl indene-1-yl lidene)methylamine-N-oxide
- 38) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate isopropyl amide
- 39) 1-(trans-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate cyclohexyl amide

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40) [1-(trans-methylimino-N-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2vl1morpholine-4-vl-methanone

- 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-yl-ethoxy)-3-phenyl-1Hindene-2-carboxylate cyclohexyl amide
- 1-(trans-methylimino-N-oxy)-3-phenyl-5-(3-phenylpropoxy)-1H-indene-2carboxvlate ethyl ester
- 43) 1-(trans-methylimino-N-oxy)-6-phenethyloxymethyl-3-phenyl-1H-indene-2-
- carboxvlate ethyl ester
- 44) (6-methoxy-3-phenylindene-1-yllidene)methylamine-N-oxide
- 1-(cis-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1Hindene-2-carboxylate ethyl ester
- 6-(2-bromoethoxy)-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-2carboxylate ethyl ester
- 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-47) indene-2-carboxylate tert-buthyl ester
- 1-(trans-methylimino-N-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-48) indene-2-carboxylate ethyl ester
- 49) 4-[2-isopropylcarbamoyl-3-(trans-methylimino-N-oxy)-1-phenyl-3H-indene-5-vi-oxylmethyl]benzoate methyl ester
- 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-50) indene-2-carboxylate isopropyl amide
- 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1Hindene-2-carboxylate cyclopropyl amide
- 3-(3-fluorophenyl)-1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-52) vlethoxy)-1H-indene-2-carboxylate isopropyl amide
- 53) (6-methoxy-1-(trans-methylimino-N-oxy)-3-phenyl-1H-indene-2-vl)acetate ethyl ester
- (6-methoxy-1-(cis-methylimino-N-oxy)-3-phenyl-1H-indene-2-yl)acetate 54) ethyl ester
- 5-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(trans-methylimino-N-oxy)-3-phenyl-55) 1H-indene-2-carboxylate isopropyl amide
- 1-(trans-methylimino-N-oxy)-6-(3-phenylpropoxy)-3-p-tolyl-1H-indene-2carboxylate ethyl ester
- 1-(trans-methylimino-N-oxy)-6-(3-phenylpropoxy)-3-thiophene-2-yl-1H-57) indene-2-carboxylate ethyl ester
- 58) 3-(4-chlorophenyl)-1-(trans-methylimino-N-oxy)-6-(3-phenylpropoxy)-1Hindene-2-carboxylate ethyl ester
- 3-(5-chlorothiophene-2-yl)-1-(trans-methylimino-N-oxy)-6-(3-phenyl 59) propoxy)-1H-indene-2-carboxylate ethyl ester
- 1-(trans-methylimino-N-oxy)-6-(3-phenylpropoxy)-3-m-tolyl-1H-indene-2carboxvlate ethyl ester

61) 1-(*trans*-methylimino-*N*-oxy)-3-(4-phenoxyphenyl)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester

- 62) 3-benzo-[1,3]-dioxol-5-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenyl propoxy)-1H-indene-2-carboxylate ethyl ester
- 63) methyl-[6-(3-phenylpropoxy)-3-pyridine-2-yl-indene-1-yllidene]-amine-*N*-oxide
- 64) 3-furan-2-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 65) 3-ethyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 66) 3-methyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 67) 1-(trans-methylimino-N-oxy)-6-(3-phenylpropoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 68) 3-cyclopropyl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 69) 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate ethyl ester
- 70) 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenyl propoxy)-1H-indene-2-carboxylate ethyl ester
- 71) 3-(1H-imidazole-4-yl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 72) 3-(1-ethyl propyl)-1-(*trans*-methylimino-*N*-oxy)-6-(3-phenylpropoxy)-1H-indene-2-carboxylate ethyl ester
- 73) 1-(trans-methylimino-N-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carboxylate amide
- 74) 6-(4-benzylmorpholine-2-ylmethoxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 75) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(3-phenylpropoxy)-1H-indene-2-carbonitrile
- 76) 1-(*trans*-methylimino-*N*-oxy)-5,6-methylenedioxy-1-oxo-3-phenyl-1H-phenyl-2-carboxylate isopropyl amide
- 77) 1-(*trans*-methylimino-*N*-oxy)-6-morpholine-4-ylmethyl-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 78) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 79) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 80) 1-(trans-methylimino-N-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 81) 6-[2-(5-ethylpyridine-2-yl)ethoxy]-1-(*trans*-methylimino-N-oxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide

82) methyl-[6-(2-morpholine-4-ylethoxy)-3-phenylindene-1-yllidene]amine-*N*-oxide

- 83) 5,6-bis-methanesulfonyloxy-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 84) 1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isobutyl ester
- 85) 1-(*trans*-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester
- 86) 1-(cis-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate methyl ester
- 87) 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate propyl ester
- 88) 3-(4-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 89) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-ylmethoxy)-1H-indene-2-carboxylate ethyl ester
- 90) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-6-(pyridine-2-yloxy)-1H-indene-2-carboxylate ethyl ester
- 91) 6-(3-methoxybenzyloxy)-1-(*trans*-methylimino-*N*-oxy)-3-phenyl-1H-indene-2-carboxylate ethyl ester
- 92) 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-thiophene-3-yl-1H-indene-2-carboxylate isopropyl amide
- 93) 3-(1-ethylpropyl)-1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 94) 3-benzo-[b]-thiophene-3-yl-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 95) 3-(4-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 96) 3-(1-ethylpropyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 97) 1-(trans-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-(2,4,6-trimethylphenyl)-1H-indene-2-carboxylate ethyl ester
- 98) 3-(2,6-dimethylphenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-morpholine-4-ylethoxy)-1H-indene-2-carboxylate ethyl ester
- 99) 1-(*trans*-methylimino-*N*-oxy)-3-phenyl-5-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide
- 100) 1-(*trans*-methylimino-*N*-oxy)-5-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl amide
- 101) 1-(cis-methylimino-N-oxy)-6-(2-morpholine-4-ylethoxy)-3-phenyl-1H-indene-2-carboxylate isopropyl ester
- 102) 3-(3-fluorophenyl)-1-(*trans*-methylimino-*N*-oxy)-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl amide

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oxy)-1H-indene-2-carboxylate ethyl ester, and

105) 1-(*trans*-methylimino-N-oxy)-3-phenyl-6-(2-pyridine-2-ylethoxy)-1H-indene-2-carboxylate isopropyl ester.

5. (currently amended) A process for preparing the indene derivative of claim 1, which comprises the step of subjecting an indenone compound of formula (II) to a condensation reaction with R₁NHOH to obtain a compound of formula (I); or comprises the steps of subjecting an indenone compound of formula (II) to a condensation reaction with NH₂OH to obtain a compound of formula (III), and conducting a reaction of the compound of formula (III) with R₁X to obtain a compound of formula (I):

$$\begin{matrix} \begin{matrix} R_6 & \begin{matrix} R_7 & \\ \\ R_8 & \begin{matrix} R_4 \end{matrix} & \begin{matrix} R_5 \end{matrix} \end{matrix} & (II)$$

wherein.

X is halogen;

cycloalkyl; and

 R_1 is C_{1-6} alkyl, C_{1-6} alkenyl, or C_{3-6} cycloalkyl, each of which is unsubstituted or substituted substituted with one or more phenyl groups;

being each unsubstituted or substitutied substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a, phenyloxy, C₁₋₆ alkyl, and C₃₋₆

 R_4 , R_5 , R_6 , and R_7 are each independently H, OH, OSO₂CH₃, O(CH₂)_mR^e, CH₂R^f, OCOCH₂OR^g, OCH₂CH₂OR^g e_f, OCH₂CH=CHR^g, or pyridine-2-yloxy, or R_5 and R_6 together form OCH₂O;

in which R^a is H, C_{1-6} alkyl, or C_{3-6} cycloalkyl, C_{1-6} alkyl and C_{3-6} cycloalkyl being each unsubstituted or substituted substituted with one or more halogens;

 R^b and R^c are each independently H, C_{1-6} alkyl, or C_{3-6} cycloalkyl; R^d is O. S. or NR^a :

COORa:

R^g is phenyl, which is unsubstituted or substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, and OR^a: and

m is an integer in the range of 1 to 5.

- 6. (currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - reacting compounds of formula (V) and (VI) to obtain a compound of formula (VII);
 - subjecting the compound of formula (VII) to cyclization to obtain a compound of formula (VIII); and
 - 3) subjecting the compound of formula (VIII) to oxidation[-],

$$\begin{array}{c} \bigcap\limits_{R_3} R_2 & (V) \\ \\ \bigcap\limits_{R_6} R_7 & Z \\ \\ \bigcap\limits_{R_6} R_7 & (VI) \\ \\ \bigcap\limits_{R_6} R_7 & (VII) \\ \\ \bigcap\limits_{R_6} R_7 & (VII) \\ \\ \bigcap\limits_{R_6} R_7 & (VIII) \\ \\ \end{array}$$
 wherein.

 $\rm R_2$ to $\rm R_7$ have the same meanings as defined in claim 5, and Z is halogen or activated leaving group.

- 7. (currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) reacting compounds of formula (IX) and (X) to obtain a compound of formula (XI):
 - 2) subjecting the compound of formula (XI) to cyclization to obtain a compound of formula (XII); and
 - 3) subjecting the compound of formula (XII) to oxidation[-],

(X)

wherein,

R₂ to R₇ have the same meanings as defined in claim 5.

- 8. (currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) reacting compounds of formula (IX) and (XIII) to obtain a compound of formula (XIV); and
 - 2) subjecting the compound of formula (XIV) to cyclization[-].

wherein.

R₂ to R₇ have the same meanings as defined in claim 5.

- 9. (currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) subjecting a compound of formula (XV) to bromination obtain a compound of formula (XVI); and
 - 2) subjecting the compound of formula (XVI) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophile[-].

wherein,

R₃ to R₇ have the same meanings as defined in claim 5.

- 10. (Currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by a process comprising the steps of:
 - 1) subjecting a compound of formula (XVII) to bromination to obtain a compound of formula (XVIII); and
 - subjecting the compound of formula (XVIII) to a carbon-carbon coupling reaction in the presence of a metal catalyst, or to a substitution reaction using a nucleophilef.

wherein.

R₂ and R₄ to R₇ have the same meanings as defined in claim 5.

11. (currently amended) The process of claim 5, wherein the indenone compound of formula (II) is prepared by subjecting a compound of formula (XIX) to an acylation reaction, a halogenation reaction followed by a substitution reaction by a nucleophile, or a carbon-carbon coupling reaction in the presence of a metal catalyst[.]

wherein.

 R_2 and R_3 have the same meanings as defined in claim 5, Y is hydroxy, thiol, amino C_{1-6} alkyl or halogen, and n is an integer in the range of 0 to 5.

12. (Currently amended) A pharmaceutical composition for modulating the activities of peroxisome proliferator activated receptors (PPARs) receptor gamma sub type comprising the compound or salt defined in claim 1 as an active ingredient together with a pharmaceutically acceptable carrier.

13. (Canceled).